

Supporting Information

Gold Decorated on 3D 2,6-Diaminopyridine Network; A Robust Catalyst for Bromination of Aromatic Compounds

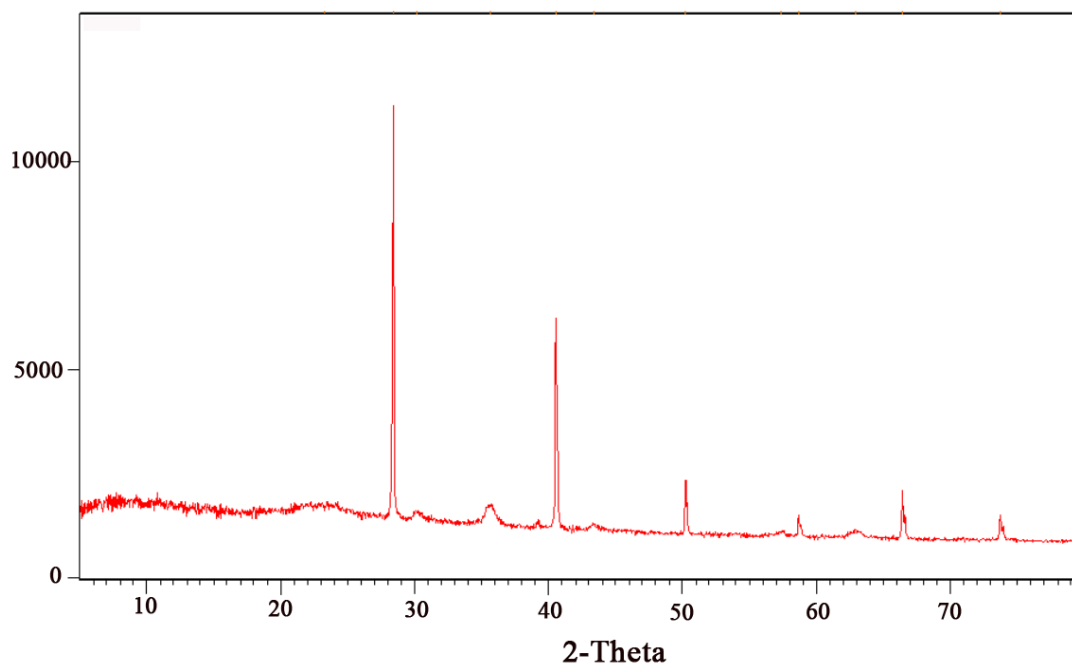
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S1. XRD pattern of MNP@ DAPN

Spectral data of the compounds:

1-bromo-4-methylbenzene:

^1H NMR (500 MHz, CDCl_3) δ 7.42 (d, J = 8.50 Hz, 2H), 7.10 (d, J = 8 Hz, 2H), 2.36 (s, 3H); ^{13}C NMR (125 MHz, CDCl_3) δ 136, 131, 130, 119, 20.9.

1-bromo-4-methoxybenzene:

^1H NMR (500 MHz, CDCl_3) δ 7.41 (d, J = 7 Hz, 2H), 6.83(d, J = 7 Hz, 2H), 3.79 (s, 3H); ^{13}C NMR (125 MHz, CDCl_3) δ 158.72, 132.25, 115.76, 112.81, 55.41.

1-bromo-4-iodobenzene:

^1H NMR (500 MHz, CDCl_3) δ 7.50 (d, J = 8.5 Hz, 2H), 7.25 (d, J = 8.25 Hz, 2H); ^{13}C NMR (125 MHz, CDCl_3) δ 139, 133, 122, 92.

4-bromoaniline:

^1H NMR (500 MHz, CDCl_3) δ 7.28(d, J = 9.5 Hz, 2H), 6.57(d, J = 8.5 Hz, 2H), 3.68(s, 2H); ^{13}C NMR (125 MHz, CDCl_3) δ 154, 132, 116, 110.

1-bromo-4-chlorobenzene:

^1H NMR (500 MHz, CDCl_3) δ 7.44 (d, J = 8.5 Hz, 2H), 7.24 (d, J = 8.5 Hz, 2H); ^{13}C NMR (125 MHz, CDCl_3) δ 133.22, 132, 130.17, 120.3.

4-bromo-1, 1'-biphenyl:

^1H NMR (500 MHz, CDCl_3) δ 7.72-7.71(m, 4H), 7.71-7.70 (m, 2H), 7.61-7.60 (m, 1H), 7.58-7.52 (m, 2H); ^{13}C NMR (125 MHz, CDCl_3) δ 140.16, 140, 131.93, 128.97, 128.8, 127.71, 127, 121.

1-bromonaphthalene:

^1H NMR (500 MHz, CDCl_3) δ 8.40(d, J = 8.5 Hz, 2H), 8.00-7.95(m, 3H), 7.78-7.69(m, 1H), 7.49-7.45(m, 1H); ^{13}C NMR (125 MHz, CDCl_3) δ 143.7, 132.1, 130, 128.4, 128.06, 127.4, 127.2, 126.8, 126.3, 122.9.

Bromobenzene:

^1H NMR (500 MHz, CDCl_3) δ 7.56-7.54 (d, J = 7.5 Hz, 2H), 7.35-7.26(m, 3H); ^{13}C NMR (125 MHz, CDCl_3) δ 132, 130.5, 127.3, 123.

9-bromophenanthrene:

^1H NMR (500 MHz, CDCl_3) δ 8.69-7.64(m, 2H), 8.41-8.39 (m, 2H), 8.12 (s, 1H), 7.82-7.61(m, 4H); ^{13}C NMR (125 MHz, CDCl_3) δ 132.2, 131.3, 130.5, 130.3, 129.7, 128.1, 127.8, 127.5, 127.4, 127.2, 127, 122.8, 122.7, 121.7.

5-bromo-2-hydroxybenzaldehyde:

^1H NMR (500 MHz, CDCl_3) δ 10.94 (s, 1H), 9.85(s, 1H), 7.68-7.50 (1H, m), 7.28 (s, 1H), 6.92(d, J = 8.5 Hz, 1H); ^{13}C NMR (125 MHz, CDCl_3) δ 195.4, 160.5, 139.6, 135.6, 121.7, 119.8, 111.3.

1-bromo-3, 5-dinitrobenzene:

^1H NMR (500 MHz, CDCl_3) δ 9.01 (s, 1H), 8.70 (s, 2H) ^{13}C NMR (125 MHz, CDCl_3) δ 117.7, 121.7, 135.6, 149.6.

5-bromoisophthalic acid:

^1H NMR (500 MHz, CDCl_3) 8.39 (s, 1H), δ 8.12 (s, 2H); ^{13}C NMR (125 MHz, $(\text{CD}_3)_2\text{SO}$) δ 165.5, 135.6, 131.9, 128.8, 121.6.

3-bromo-5-nitrosalicylaldehyde:

^1H NMR (500 MHz, CDCl_3) δ 12.24 (s, 1H), 9.85 (s, 1H), 8.68 (s, 1H), 8.48 (s, 1H); ^{13}C NMR (125 MHz, CDCl_3) δ 195.4, 160.5, 140.1, 135.6, 128.8, 119.8, 111.3.

3-bromo-5-nitrobenzoic acid:

^1H NMR (500 MHz, CDCl_3) δ 8.96 (s, 1H), 8.64 (s, 1H), 8.45 (s, 1H); ^{13}C NMR (125 MHz, $(\text{CD}_3)_2\text{SO}$) δ 164.5, 148.6, 135.6, 132.2, 129.7, 121.7, 119.8.

4-bromophenol:

^1H NMR (500 MHz, CDCl_3) 7.82 (d, J = 7.5 Hz, 2H), δ 7.57 (d, J = 7.5 Hz, 2H), 5.12 (s, 1H); ^{13}C NMR (125 MHz, CDCl_3) δ 158.7, 132.6, 121.6, 116.7.

1-bromo-4-fluorobenzene:

^1H NMR (500 MHz, CDCl_3) δ 7.28 (d, J = 9.5 Hz, 2H), 7.69 (d, J = 7 Hz, 2H); ^{13}C NMR (125 MHz, CDCl_3) δ 160.1, 135.6, 134.7, 121.7.

1-iodo-3, 5-dinitrobenzene:

^1H NMR (500 MHz, CDCl_3) δ 8.72 (s, 1H), 8.08 (s, 2H) ^{13}C NMR (125 MHz, CDCl_3) δ 93.5, 122.1, 134.4, 143.5.

1-chloro-3, 5-dinitrobenzene:

^1H NMR (500 MHz, CDCl_3) δ 8.64 (s, 1H), 8.86 (s, 2H) ^{13}C NMR (125 MHz, CDCl_3) δ 116.7, 122.1, 125.5, 143.3.